

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP120520\
 Data File : BP004165.D
 Acq On : 05 Dec 2020 17:04
 Operator : CG/JU
 Sample : PB133233BL
 Misc : MDL-WATER-BL-03
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK233

Quant Time: Dec 07 03:02:28 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP120420.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Dec 07 01:39:51 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.85	152	377119	20.00	ng/ul	0.00
20) Naphthalene-d8	10.65	136	1549150	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.49	164	972283	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.25	188	2141657	20.00	ng/ul	0.00
79) Chrysene-d12	21.35	240	2261824	20.00	ng/ul	0.00
88) Perylene-d12	23.71	264	2457179	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.34	96	59973	6.11	ng/uL	0.00
4) Pyridine-d5	3.74	84	842955	30.23	ng/ul	0.00
7) Phenol-d5	7.00	99	1004687	29.80	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.18	67	621576	31.89	ng/ul	0.00
11) 2-Chlorophenol-d4	7.38	132	775271	29.55	ng/ul	0.00
15) 4-Methylphenol-d8	8.55	113	759292	28.59	ng/ul	0.00
21) Nitrobenzene-d5	9.00	128	399649	31.64	ng/ul	0.00
24) 2-Nitrophenol-d4	9.72	143	340204	29.50	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.26	165	709403	27.95	ng/ul	0.00
31) 4-Chloroaniline-d4	10.78	131	1097772	29.48	ng/ul	0.00
46) Dimethylphthalate-d6	13.90	166	2502075	32.40	ng/ul	0.00
49) Acenaphthylene-d8	14.19	160	2856321	30.10	ng/ul	0.00
54) 4-Nitrophenol-d4	14.67	143	344356	27.25	ng/ul	0.00
60) Fluorene-d10	15.49	176	2111715	31.27	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.60	200	212448	24.41	ng/ul	0.00
73) Anthracene-d10	17.35	188	3365903	31.98	ng/ul	0.00
81) Pyrene-d10	19.59	212	3939922	33.11	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.56	264	4413531	33.44	ng/ul	0.00

Target Compounds Ovalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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